





## 10522150.tsm



[illegible][illegible][illegible][illegible]

## 10522150.trn

[illegible][illegible][illegible][illegible]

## 10522150.trn

[illegible][illegible]

**NAME** 3-(4-*tert*-butylphenyl)-2-phenyl-5-thiothiazole  
**SMILES** CC(C)(C)c1ccc(cc1)-c2cc3cc(sc3cc2-c4ccccc4)S  
**MO** 157470-00-0 CAS#157470  
**DE** 1-Methoxy-2-phenyl-3-(4-*tert*-butylphenyl)-5-thiothiazole  
**SMILES** COc1cc2cc3cc(sc3cc2-c4ccccc4)-c5ccc(cc5)C(C)(C)C  
**Abbildung**   
**NAME** 3-(4-*tert*-butylphenyl)-2-phenyl-5-thiothiazole  
**SMILES** CC(C)(C)c1ccc(cc1)-c2cc3cc(sc3cc2-c4ccccc4)S  
**MO** 157470-70-2 CAS#157470  
**DE** 1-Methoxy-2-phenyl-3-(4-*tert*-butylphenyl)-5-thiothiazole  
**SMILES** COc1cc2cc3cc(sc3cc2-c4ccccc4)-c5ccc(cc5)C(C)(C)C  
**Abbildung** 

[illegible]

## 10522150.trn

[illegible][illegible][illegible][illegible]

```

uu 015730-01-8 EMPL01
cu ?=ahshaylo{,},?|root me, ?={0-1|notky|-}0+intl+0+||+>
pyridadshy|ovj^, {0)+ CH 10000 00000

```

[illegible]O=C1C=CC(=C2C=CC(=C3C=CC(=C2)C=C3)C=C1)C4=CC=CC=C4


CRY 3  
 FROM 316-37-2  
 CRY 74 86 86

O=C1C=CC(=C1)C2=CC=CC=C2C3=CC=CC=C3[illegible]c1ccc2c(c1)c(c3ccccc23)C(=O)O

DN 3  
CNR 65281-66-6  
CNR C19 628 04 8

```
Ans 016730-7)-S ChR19N
Cu 1-16d 016730-7)-S ChR19N
1101-J-16m 016730-7)-S ChR19N
(ChR 16m 016730-7)-S ChR19N
(ChR 16m 016730-7)-S ChR19N
(ChR 16m 016730-7)-S ChR19N
```

**Abstract:** **SYNTHESIS OF**



The chemical structure shows a central benzene ring substituted at the 1 and 5 positions. At the 1-position, there is a pyrrolidine ring attached via its nitrogen atom. At the 5-position, there is a 1,2,3,4-tetrahydronaphthalen-1-yl group attached via its 1-position. The tetrahydronaphthalene system consists of a benzene ring fused to a cyclohexane ring, with the attachment point at the 1-position of the cyclohexane ring.

CIV 1  
Q No 76-01-1  
100 C1 M P3 G3

Chem: bond geometry as above

RD 00167-75-3 CAPLON

CD 1-ethynyl-2-methyl-2-butene,  
trans-2,3,4-trimethylpent-2-ene

ON 3  
 EMB 855758-45-9  
 TWE C19 020 04 01

CE 1-(4-chlorophenyl)-3,5-diisopropyl-2-[(5-{2-methyl-5-oxooctan-2-yl}-2-pyrimidinyl)oxy]-4-nitrobenzylidene (11), (12):  $[\alpha]_D^{20} + 100.00$  (CHCl<sub>3</sub>)

Isomeric stereoisomerism.

Chemical structure of compound 11: CC(C)C1=CC=C(C=C1C2=CC=CC=C2C3=CC(=CC=C3)C(=O)N(C)C)C4=CC=C(C=C4)C5=CC=CC=C5C6=CC=CC=C6C7=CC=CC=C7C8=CC=CC=C8C9=CC=CC=C9C10=CC=CC=C10C11=CC=CC=C11C12=CC=CC=C12C13=CC=CC=C13C14=CC=CC=C14C15=CC=CC=C15C16=CC=CC=C16C17=CC=CC=C17C18=CC=CC=C18C19=CC=CC=C19C20=CC=CC=C20C21=CC=CC=C21C22=CC=CC=C22C23=CC=CC=C23C24=CC=CC=C24C25=CC=CC=C25C26=CC=CC=C26C27=CC=CC=C27C28=CC=CC=C28C29=CC=CC=C29C30=CC=CC=C30C31=CC=CC=C31C32=CC=CC=C32C33=CC=CC=C33C34=CC=CC=C34C35=CC=CC=C35C36=CC=CC=C36C37=CC=CC=C37C38=CC=CC=C38C39=CC=CC=C39C40=CC=CC=C40C41=CC=CC=C41C42=CC=CC=C42C43=CC=CC=C43C44=CC=CC=C44C45=CC=CC=C45C46=CC=CC=C46C47=CC=CC=C47C48=CC=CC=C48C49=CC=CC=C49C50=CC=CC=C50C51=CC=CC=C51C52=CC=CC=C52C53=CC=CC=C53C54=CC=CC=C54C55=CC=CC=C55C56=CC=CC=C56C57=CC=CC=C57C58=CC=CC=C58C59=CC=CC=C59C60=CC=CC=C60C61=CC=CC=C61C62=CC=CC=C62C63=CC=CC=C63C64=CC=CC=C64C65=CC=CC=C65C66=CC=CC=C66C67=CC=CC=C67C68=CC=CC=C68C69=CC=CC=C69C70=CC=CC=C70C71=CC=CC=C71C72=CC=CC=C72C73=CC=CC=C73C74=CC=CC=C74C75=CC=CC=C75C76=CC=CC=C76C77=CC=CC=C77C78=CC=CC=C78C79=CC=CC=C79C80=CC=CC=C80C81=CC=CC=C81C82=CC=CC=C82C83=CC=CC=C83C84=CC=CC=C84C85=CC=CC=C85C86=CC=CC=C86C87=CC=CC=C87C88=CC=CC=C88C89=CC=CC=C89C90=CC=CC=C90C91=CC=CC=C91C92=CC=CC=C92C93=CC=CC=C93C94=CC=CC=C94C95=CC=CC=C95C96=CC=CC=C96C97=CC=CC=C97C98=CC=CC=C98C99=CC=CC=C99C100=CC=CC=C100C101=CC=CC=C101C102=CC=CC=C102C103=CC=CC=C103C104=CC=CC=C104C105=CC=CC=C105C106=CC=CC=C106C107=CC=CC=C107C108=CC=CC=C108C109=CC=CC=C109C110=CC=CC=C110C111=CC=CC=C111C112=CC=CC=C112C113=CC=CC=C113C114=CC=CC=C114C115=CC=CC=C115C116=CC=CC=C116C117=CC=CC=C117C118=CC=CC=C118C119=CC=CC=C119C120=CC=CC=C120C121=CC=CC=C121C122=CC=CC=C122C123=CC=CC=C123C124=CC=CC=C124C125=CC=CC=C125C126=CC=CC=C126C127=CC=CC=C127C128=CC=CC=C128C129=CC=CC=C129C130=CC=CC=C130C131=CC=CC=C131C132=CC=CC=C132C133=CC=CC=C133C134=CC=CC=C134C135=CC=CC=C135C136=CC=CC=C136C137=CC=CC=C137C138=CC=CC=C138C139=CC=CC=C139C140=CC=CC=C140C141=CC=CC=C141C142=CC=CC=C142C143=CC=CC=C143C144=CC=CC=C144C145=CC=CC=C145C146=CC=CC=C146C147=CC=CC=C147C148=CC=CC=C148C149=CC=CC=C149C150=CC=CC=C150C151=CC=CC=C151C152=CC=CC=C152C153=CC=CC=C153C154=CC=CC=C154C155=CC=CC=C155C156=CC=CC=C156C157=CC=CC=C157C158=CC=CC=C158C159=CC=CC=C159C160=CC=CC=C160C161=CC=CC=C161C162=CC=CC=C162C163=CC=CC=C163C164=CC=CC=C164C165=CC=CC=C165C166=CC=CC=C166C167=CC=CC=C167C168=CC=CC=C168C169=CC=CC=C169C170=CC=CC=C170C171=CC=CC=C171C172=CC=CC=C172C173=CC=CC=C173C174=CC=CC=C174C175=CC=CC=C175C176=CC=CC=C176C177=CC=CC=C177C178=CC=CC=C178C179=CC=CC=C179C180=CC=CC=C180C181=CC=CC=C181C182=CC=CC=C182C183=CC=CC=C183C184=CC=CC=C184C185=CC=CC=C185C186=CC=CC=C186C187=CC=CC=C187C188=CC=CC=C188C189=CC=CC=C189C190=CC=CC=C190C191=CC=CC=C191C192=CC=CC=C192C193=CC=CC=C193C194=CC=CC=C194C195=CC=CC=C195C196=CC=CC=C196C197=CC=CC=C197C198=CC=CC=C198C199=CC=CC=C199C200=CC=CC=C200C201=CC=CC=C201C202=CC=CC=C202C203=CC=CC=C203C204=CC=CC=C204C205=CC=CC=C205C206=CC=CC=C206C207=CC=CC=C207C208=CC=CC=C208C209=CC=CC=C209C210=CC=CC=C210C211=CC=CC=C211C212=CC=CC=C212C213=CC=CC=C213C214=CC=CC=C214C215=CC=CC=C215C216=CC=CC=C216C217=CC=CC=C217C218=CC=CC=C218C219=CC=CC=C219C220=CC=CC=C220C221=CC=CC=C221C222=CC=CC=C222C223=CC=CC=C223C224=CC=CC=C224C225=CC=CC=C225C226=CC=CC=C226C227=CC=CC=C227C228=CC=CC=C228C229=CC=CC=C229C230=CC=CC=C230C231=CC=CC=C231C232=CC=CC=C232C233=CC=CC=C233C234=CC=CC=C234C235=CC=CC=C235C236=CC=CC=C236C237=CC=CC=C237C238=CC=CC=C238C239=CC=CC=C239C240=CC=CC=C240C241=CC=CC=C241C242=CC=CC=C242C243=CC=CC=C243C244=CC=CC=C244C245=CC=CC=C245C246=CC=CC=C246C247=CC=CC=C247C248=CC=CC=C248C249=CC=CC=C249C250=CC=CC=C250C251=CC=CC=C251C252=CC=CC=C252C253=CC=CC=C253C254=CC=CC=C254C255=CC=CC=C255C256=CC=CC=C256C257=CC=CC=C257C258=CC=CC=C258C259=CC=CC=C259C260=CC=CC=C260C261=CC=CC=C261C262=CC=CC=C262C263=CC=CC=C263C264=CC=CC=C264C265=CC=CC=C265C266=CC=CC=C266C267=CC=CC=C267C268=CC=CC=C268C269=CC=CC=C269C270=CC=CC=C270C271=CC=CC=C271C272=CC=CC=C272C273=CC=CC=C273C274=CC=CC=C274C275=CC=CC=C275C276=CC=CC=C276C277=CC=CC=C277C278=CC=CC=C278C279=CC=CC=C279C280=CC=CC=C280C281=CC=CC=C281C282=CC=CC=C282C283=CC=CC=C283C284=CC=CC=C284C285=CC=CC=C285C286=CC=CC=C286C287=CC=CC=C287C288=CC=CC=C288C289=CC=CC=C289C290=CC=CC=C290C291=CC=CC=C291C292=CC=CC=C292C293=CC=CC=C293C294=CC=CC=C294C295=CC=CC=C295C296=CC=CC=C296C297=CC=CC=C297C298=CC=CC=C298C299=CC=CC=C299C300=CC=CC=C300C301=CC=CC=C301C302=CC=CC=C302C303=CC=CC=C303C304=CC=CC=C304C305=CC=CC=C305C306=CC=CC=C306C307=CC=CC=C307C308=CC=CC=C308C309=CC=CC=C309C310=CC=CC=C310C311=CC=CC=C311C312=CC=CC=C312C313=CC=CC=C313C314=CC=CC=C314C315=CC=CC=C315C316=CC=CC=C316C317=CC=CC=C317C318=CC=CC=C318C319=CC=CC=C319C320=CC=CC=C320C321=CC=CC=C321C322=CC=CC=C322C323=CC=CC=C323C324=CC=CC=C324C325=CC=CC=C325C326=CC=CC=C326C327=CC=CC=C327C328=CC=CC=C328C329=CC=CC=C329C330=CC=CC=C330C331=CC=CC=C331C332=CC=CC=C332C333=CC=CC=C

④ NCJ

NO 006197-00-0 CNF/OW  
CR I-S&MILPUL01.1.1 004 AN+. P-[19-(1-phonyl)-5-hydroxymethyl]-  
pyrimidinyl[carb], hydrochloride (1:1), (10)\* NCH [UNCL CLASS]

```

      = FM      1
      CDR      215130-50-2
      CDR      C18 019 05 01
      DD-00107  01-7-0000-00.07.72.

```

FM	3
FAM	70-93-3
EW	03 06 08 02

c1ccc2c(c1)c(c3ccccc23)C(=O)NCC4CCCC4

500 3  
 500 110-117-4  
 500 110 117 4

CC1=CC=CC=C1C(=O)O  
 SMILES: CC1=CC=CC=C1C(=O)O

10

17	003101-70-20	015201-50-00	055201-67-10
	055202-71-00	015700-00-00	055710-67-70
	105710-00-00	015700-01-20	105710-00-00
	105720-01-10	055700-01-70	055710-72-00

[illegible][illegible]


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NO 215736-01-2  CML00
CV  COUNCILMAN, B-19-11781-2-AMMIGUEL012.2.1JACK-B-GARY-1-
    PPR00000012-2-AMMIGUEL1, 1.1-GARY-1-JACK-AMMIGUEL
Absolute chronometry.

```

[illegible]

As values of  $\alpha$  increase, the


  
**IN** 054719-00-0 CHEMREX
   
**CU** 1-(4-bromophenyl)-2,3-dihydro-1H-indene 7-((E)-2-(4-bromophenyl)-1-propenyl)-2,3-dihydro-1H-indene

4466 J. Neurosci., June 1, 1997, 17(12):4462-4466




NO 055723-49-9 DCE

Chemical structure of 2,6-dimethylphenol (Cresol): Cc1cc(C)cc(O)c1

  
In 155710-65-7 [26]60  
FE Acetamide, O-[2-(4-(4-chlorophenyl)-2-propenyl)-p-toluidinyl]phenyl[<sup>a</sup>] ccr 10000 (new)

[illegible]


  
 SMILES: CC1=C(C)C(=O)N(C1)C
  
 IUPAC: 2,4,6-trimethyl-3-pyridinol

































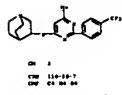


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## 10522150.trn

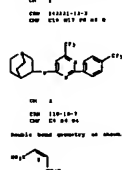
© 2006 Blackwell Publishing Ltd  
Journal compilation © 2006 Blackwell Publishing Ltd



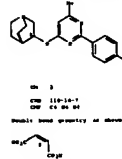
```

IN 141221-13-4 CMLFV9
CE 2-Acetylcholinesterase, 2-[16-(4-fluoromethyl)-2-(4-
   [4-fluoromethyl]phenyl)-4-pyridylidene]pyrrolidine-1-oxide-1-oxide
INCL 100 1000 0001

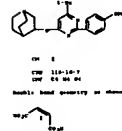
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14. NUMBER OF ACRES OF FORESTED LAND 1947 529 88 273 150,000

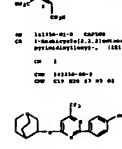


CM 1  
 CMM 141441-20-3  
 CMT C22 R39 03 03

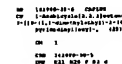


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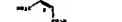
14. NUMBER OF SETS OF CHANGES CONTAINED: 207 SETS OF 225 (207/225=92%)



Chemical structure of 1,2-dichloroethane (DCE): ClCCl



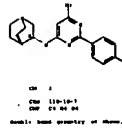
14 JANUARY 1987 10 02/50 COPY/1007 1007 ACS on 978 100001000000



Page 235

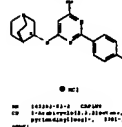
## 10522150.trn

1-6. APPROVED 28 MAY 58. CALIFORNIA COPYRIGHT 1957 ACE AND STE. HANFORD, CALIF.

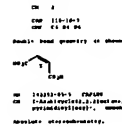
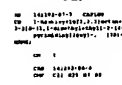


```

ME 10JIV-00-0 CBJ/BO
FO [-Aambirysia[4.8]ortuu, 1-312-(4-fiwerephayil-6-watig[-o-
pywimayillori-, ammbiporoph[ortu, (A) (-WZ) ICH 1994E MMBZ
MashGt: Microthermistry.
```



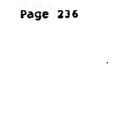
14 NUMBER 13 OF 38 CHANGES EMPLOYMENT 1997 Q4 vs 97Q3 (THOUSANDS)

ClC1=CC=CC=C1

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## 10522150.trn

14. APPROX 10:00:30 EARLIER. COPYRIGHT 2005 BOWEN LTD. (Don't Lapse)



=> FIL STNGUIDE  
 FILE 'STNGUIDE' ENTERED AT 13:27:55 ON 05 NOV 2007  
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FILE CONTAINS CURRENT INFORMATION.  
LAST RELOADED: Nov 2, 2007 (20071102/UP).

--> file caplus  
FILE 'CAPLUS' ENTERED AT 13:32:27 ON 05 NOV 2007  
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FILE COVERS 1907 - 5 Nov 2007 VOL 147 ISS 20  
FILE LAST UPDATED: 4 Nov 2007 (20071104/ED)

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They are available for your review at:

<http://www.cas.org/infopolicy.html>

and his

(FILE 'HOMB' ENTERED AT 13:21:34 ON 05 NOV 2007)

```

FILE 'REGISTRY' ENTERED AT 13:21:51 ON 05 NOV 2007
L1      STRUCTURE UPLOADED
L2      2 S L1
L3      449 S L1 FULL

```

FILE 'CAPLUS' ENTERED AT 13:26:23 ON 05 NOV 2007  
L4 30 S L3

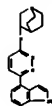
FILE 'STNGUIDE' ENTERED AT 13:27:55 ON 05 NOV 2007


FILE 'CAPLUS' ENTERED AT 13:32:27 ON 05 NOV 2007

```

=> d 12
YOU HAVE REQUESTED DATA FROM FILE 'REGISTRY' - CONTINUE? (Y)/N:Y

```

[illegible]

CN 2  
 CNB 110-17-8  
 CNF 24 00 34  
 Doubly bond geometry as shown  


2 AFFIDAVITS RE FILE CA 11497 VD DATE:  
1 AFFIDAVITS RE FILE CNLPG 11497 VD DATE:

```

=> d 12/read
'L2' MUST END IN '/Q', '/A', '/L', '/S' OR '/B'
The saved name for a query (or structure or screen set) must end with
'/Q'. The saved name for an answer set must end with '/A'. The saved
name for an L2 list must end with '/L'. SDI request names must end
with '/S'. To see a list of all saved query, answer set, and L2 list
names for this loginid, enter "DISPLAY SAVED" at an arrow
prompt (=>). Enter "DISPLAY SAVED/S" to see a list of SDI request
names. Enter "DISPLAY SAVED/B" to see a list of BATCH search
names.
ESBUSE1

```

=> FIL REGISTRY  
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DICTIONARY FILE UPDATES: 4 NOV 2007 HIGHEST RN 952404-20-5

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Please note that search-term pricing does apply when conducting SmartSELECT searches.

REGISTRY includes numerically searchable data for experimental and predicted properties as well as tags indicating availability of experimental property data in the original document. For information on property searching in REGISTRY, refer to:

<http://www.cag.org/support/stngen/stndoc/properties.html>

21 per d

[illegible]

1 2 3 4 5





## EAST Search History

Ref #	Hits	Search Query	DBs	Default Operator	Plurals	Time Stamp
L1	666	514/305.ccls.	US-PGPUB; USPAT; USOCR; FPRS; EPO; JPO; DERWENT; IBM_TDB	OR	ON	2007/11/05 14:44
L2	159	L1 and muscarinic	US-PGPUB; USPAT; USOCR; FPRS; EPO; JPO; DERWENT; IBM_TDB	OR	ON	2007/11/05 14:45
L3	9	"5512574"	US-PGPUB; USPAT; USOCR; FPRS; EPO; JPO; DERWENT; IBM_TDB	OR	ON	2007/11/05 14:46
L4	4	"5888999"	US-PGPUB; USPAT; USOCR; FPRS; EPO; JPO; DERWENT; IBM_TDB	OR	ON	2007/11/05 14:46
L5	47	"04333"	US-PGPUB; USPAT; USOCR; FPRS; EPO; JPO; DERWENT; IBM_TDB	OR	ON	2007/11/05 14:48
L6	0	"04333 A1"	US-PGPUB; USPAT; USOCR; FPRS; EPO; JPO; DERWENT; IBM_TDB	OR	ON	2007/11/05 14:49
L7	1241	chokai.in.	US-PGPUB; USPAT; USOCR; FPRS; EPO; JPO; DERWENT; IBM_TDB	OR	ON	2007/11/05 14:49

## EAST Search History

L8	6	L7 and pyrimidine	US-PGPUB; USPAT; USOCR; FPRS; EPO; JPO; DERWENT; IBM_TDB	OR	ON	2007/11/05 14:51
L9	372	"12711"	US-PGPUB; USPAT; USOCR; FPRS; EPO; JPO; DERWENT; IBM_TDB	OR	ON	2007/11/05 14:51
L10	26601	L9 pyrazine	US-PGPUB; USPAT; USOCR; FPRS; EPO; JPO; DERWENT; IBM_TDB	OR	ON	2007/11/05 14:51
L11	29	L9 and pyrazine	US-PGPUB; USPAT; USOCR; FPRS; EPO; JPO; DERWENT; IBM_TDB	OR	ON	2007/11/05 15:17
L12	210	"20819"	US-PGPUB; USPAT; USOCR; FPRS; EPO; JPO; DERWENT; IBM_TDB	OR	ON	2007/11/05 15:17